Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	175	560/188.ccls.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L2	5626	musk	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L3	0	L1 and L2	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L4	26151	terpene	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L5	366	560/252.ccls.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L6	536	L1 or L5	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L7	0	L2 and L6	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L8	3	L4 and L6	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L9	53861	fragr\$	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L10	217	512/22.ccls.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L11	136	512/18.ccls.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36

L12	348	L11 or L10	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L13	115	3,3-dimethylcyclohexane	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:51
L14	9	3,3-dimethyl-cyclohexane	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L15	1	L12 and L13	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L16	10	L14 or L15	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L17	123	L14 or L13	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L18	347	3,3-dimethylcyclohex\$	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L19	92	L9 and L18	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L20	175	560/188.ccls.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L21	5626	musk	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
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L25	536	L1 or L5	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L26	0	L2 and L6	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L27	3	L4 and L6	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
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L35	92	L9 and L18	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36

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L38	1	"4534891".PN.	USPAT; USOCR	OR	ON	2005/12/29 06:36
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L49	2	("6384269").URPN.	USPAT	OR	ON	2005/12/29 06:36
L50	2	("2004234568").PN.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2005/12/29 06:36

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L59	9	3,3-dimethyl-cyclohexane	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
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L61	53	L2 and L19	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L62	11	cyclohexylalkanol	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:53
L63	22	3,3-dimethylcyclohex\$.clm.	US-PGPUB	OR	ON	2005/12/29 06:36

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L68	2	("2004234568").PN.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR .	OFF	2005/12/29 06:36
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L71	18	"2008167"	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L72	2	"3487102".PN.	USPAT; USOCR	OR	ON	2005/12/29 06:36
L73	9	"4504412"	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L74	3	L1 and L4	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L75	5	L6 and L9	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L76	5	L11 and L10	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L77	9	3,3-dimethyl-cyclohexane	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36

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L78	20	L9 and L17	US-PGPUB; USPAT; EPO; JPO;	OR .	ON	2005/12/29 06:36
L79	53	L2 and L19	DERWENT US-PGPUB; USPAT;	OR	ON	2005/12/29 06:36
L80	11	cyclohexylalkanol	EPO; JPO; DERWENT US-PGPUB;	OR	ON	2005/12/29 06:36
			USPAT; EPO; JPO; DERWENT			
L81	22	3,3-dimethylcyclohex\$.clm.	US-PGPUB	OR	ON	2005/12/29 06:36
L82	32	"4622221"	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
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L84	105	L2 and L12	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L85	136	512/18.ccls.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
L86	105	L2 and L12	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:36
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L88	0	I13 and I87	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:51
L89	138298	cyclohexyl	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 07:21
L90	27	187 and 189	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2005/12/29 06:54

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        DEC 21 IPC search and display fields enhanced in CA/CAplus with the
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                IPC reform
        DEC 23 New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/USPAT2
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```
=> e 1-Propanol, 2-methyl-2-[1-(3-methylcyclohexyl)ethoxy]-, acetate/cn
                   1-PROPANOL, 2-METHYL-2,3-DI-2-PYRIDYL-, DIPICRATE/CN
E1
                   1-PROPANOL, 2-METHYL-2, 3-DIPHENYL-/CN
E2
             0 --> 1-PROPANOL, 2-METHYL-2-1-(3-METHYLCYCLOHEXYL)ETHOXY-, ACET
E3
                   ATE/CN
                   1-PROPANOL, 2-METHYL-2-(((1,2,3,4-TETRAHYDRO-8-QUINOLINYL)ME
E4
             1
                   THYL) AMINO) -/CN
                   1-PROPANOL, 2-METHYL-2-(((1-(2-(METHYLAMINO)-4-(PROPYLAMINO)
E5
             1
                   -7-QUINAZOLINYL) -1H-PYRROL-3-YL) METHYL) AMINO) -/CN
                   1-PROPANOL, 2-METHYL-2-(((1-METHYL-1H-PYRROL-2-YL)METHYL)(PH
             1
E6
                   ENYLMETHYL) AMINO) -/CN
                   1-PROPANOL, 2-METHYL-2-(((2-(2-PROPENYL)PHENYL)METHYL)AMINO)
E7
             1
                   -/CN
                   1-PROPANOL, 2-METHYL-2-(((2-(2-PROPENYLOXY)-1-NAPHTHALENYL)M
E8
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                   ETHYL) AMINO) -/CN
                   1-PROPANOL, 2-METHYL-2-(((2-(2-PROPENYLOXY)PHENYL)METHYL)AMI
E9
             1
                   NO) -/CN
                   1-PROPANOL, 2-METHYL-2-(((2-(METHYL-D3)PHENYL)METHYL)AMINO)-
E10
             1
                   /CN
             1
                   1-PROPANOL, 2-METHYL-2-(((2-(PHENYLMETHOXY)-1-NAPHTHALENYL)M
E11
                   ETHYL) AMINO) -/CN
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E12
             1
                   NO) -/CN
```

```
=> e 1-Propanol, 2-methyl-2-(1-(3-methylcyclohexyl)ethoxy)-, acetate/cn
                   1-PROPANOL, 2-METHYL-2-(1-(2-METHYLCYCLOHEXYL)ETHOXY)-/CN
E1
                   1-PROPANOL, 2-METHYL-2-(1-(2-METHYLCYCLOHEXYL)ETHOXY)-, PROP
E2
                   ANOATE/CN
             1 --> 1-PROPANOL, 2-METHYL-2-(1-(3-METHYLCYCLOHEXYL)ETHOXY)-, ACET
E3
                   ATE/CN
                   1-PROPANOL, 2-METHYL-2-(1-(3-METHYLCYCLOHEXYL)ETHOXY)-, PROP
E4
             1
                   ANOATE/CN
             1
                   1-PROPANOL, 2-METHYL-2-(1-(5-METHYL-3-CYCLOHEXEN-1-YL)ETHOXY
E5
                   ) -/CN
                   1-PROPANOL, 2-METHYL-2-(1-(5-METHYL-3-CYCLOHEXEN-1-YL)ETHOXY
             1
E6
                   )-, PROPANOATE/CN
                   1-PROPANOL, 2-METHYL-2-(1-METHYLHEPTYLAMINO)-, BENZOATE, HYD
             1
E7
                   ROCHLORIDE/CN
                   1-PROPANOL, 2-METHYL-2-(1-METHYLHEPTYLAMINO)-, FORMATE/CN
E8
             1
                   1-PROPANOL, 2-METHYL-2-(1-METHYLHEXYLAMINO)-/CN
E9
             1
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E10
             1
                   1-PROPANOL, 2-METHYL-2-(2,4,6-TRINITROANILINO)-/CN
             1
E11
                   1-PROPANOL, 2-METHYL-2-(2-METHYL-2-((2-METHYL-2-PROPENYL)OXY
             1
E12
                   ) PROPOXY) -/CN
=> e3
             1 "1-PROPANOL, 2-METHYL-2-(1-(3-METHYLCYCLOHEXYL)ETHOXY)-, ACETATE
Ll
               "/CN
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     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
L1
     610769-93-2 REGISTRY
RN
     Entered STN: 30 Oct 2003
ED
     1-Propanol, 2-methyl-2-[1-(3-methylcyclohexyl)ethoxy]-, acetate
CN
     (9CI) (CA INDEX NAME)
     3D CONCORD
FS
     C15 H28 O3
MF
SR
     CA
                  CA, CAPLUS, CASREACT, USPATFULL
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1 L1 L2

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- Methods for the production of novel alicyclic esters having a musky smell ΤI
- 2003:796643 CAPLUS AN
- 139:307907 DN
- Methods for the production of novel alicyclic esters having a musky smell ΤI
- IN Eh, Marcus
- Symrise GmbH & Co. KG, Germany PA
- SO PCT Int. Appl., 45 pp.

CODEN: PIXXD2

- DT Patent
- LA German

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 R^6
 R^7

The invention relates to novel alicyclic esters I [R1 = Me; R2, R4 = H; R3 AB = H, Me; R5, R6 = H, Me; Y = CR7R8OC(:O)R9; R7, R8 = H, Me; R9 = C1-5-alkyl, C2-5-alkylene; or R1, R2 = Me, Et; R3, R4 = H, Me; R5R6 = O; Y = CR7R8OC(:0)R9; or R1, R2 = Me, Et; R4, R5, R6, R7 = H, Me; Y = CR7R8OC(:0)R9], methods for their production, for their use as odorous substances for perfumed products and for odorous substance mixts. containing the inventive compds. The procedure for the preparation of I is characterized by reaction of cyclohexylalkanols II with carboxylic acids [R9CO2CR7R8CO2H, R9CO2H or XCR7R8CO2H (X = OH, halogen)] anhydrides [(R9CO2)20 or (XCR7R8CO2)20], or epoxides, III. Thus, I [R1 = Me, R2 - R4] = H, R6 = CHMe2, Y = O2CEt] was prepared from 1-cyclohexylethanol via reaction with isobutylene oxide in cyclohexane containing BF3·OEt2, followed by reaction with (EtCO2)20 containing Et3N in the presence of catalytic DMAP. The odor of I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2CEt] was characterized (perceptible rose bloom note).

=> logoff hold COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.10	11.04
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.73	-0.73

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 07:17:59 ON 29 DEC 2005 Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 07:18:58 ON 29 DEC 2005 FILE 'CAPLUS' ENTERED AT 07:18:58 ON 29 DEC 2005 COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	3.10	11.04
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.73	-0.73
=> file reg	atvan nii	moma i
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	3.55	11.49
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.73	-0.73

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 28 DEC 2005 HIGHEST RN 870751-96-5 DICTIONARY FILE UPDATES: 28 DEC 2005 HIGHEST RN 870751-96-5

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search iteration limits have been increased. See HELP SLIMITS for details.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

```
=> e Acetic acid, (1-oxopropoxy)-, 1-(3,3-dimethylcyclohexyl)-1-methylethyl/cn
                   ACETIC ACID, (1-OXOPROPOXY)-, 1,2,4-TRIMETHYL-2-PENTENYL EST
E1
                   ER/CN
                   ACETIC ACID, (1-OXOPROPOXY)-, 1-(3,3-DIMETHYL-1-CYCLOHEXEN-1
E2
                   -YL) ETHYL ESTER/CN
             0 --> ACETIC ACID, (1-OXOPROPOXY)-, 1-(3,3-DIMETHYLCYCLOHEXYL)-1-M
E3
                   ETHYLETHYL/CN
                   ACETIC ACID, (1-OXOPROPOXY)-, 1-(3,3-DIMETHYLCYCLOHEXYL)-1-M
             1
E4
                   ETHYLETHYL ESTER/CN
                   ACETIC ACID, (1-OXOPROPOXY) -, 1-(3,3-DIMETHYLCYCLOHEXYL) ETHY
             1
E5
                   L ESTER/CN
                   ACETIC ACID, (1-OXOPROPOXY)-, 1-(5,5-DIMETHYL-1-CYCLOHEXEN-1
             1
E6
                   -YL) ETHYL ESTER/CN
                   ACETIC ACID, (1-OXOPROPOXY)-, 1-CYCLOHEXYL-1-METHYLETHYL EST
F.7
             1
                   ER/CN
                   ACETIC ACID, (1-OXOPROPOXY) -, 1-METHYLPROPYL ESTER/CN
             1
E8
                   ACETIC ACID, (1-OXOPROPOXY)-, 2-AMINO-2-METHYLPROPYL ESTER/C
             1
E9
                   ACETIC ACID, (1-OXOPROPOXY)-, 2-AMINO-2-METHYLPROPYL ESTER,
             1
E10
                   HYDROCHLORIDE/CN
                   ACETIC ACID, (1-OXOPROPOXY)-, 2-METHOXY-2-OXOETHYL ESTER/CN
E11
             1
                   ACETIC ACID, (1-OXOPROPOXY) -, 2-PHENOXYETHYL ESTER/CN
             1
E12
=> e4
             1 "ACETIC ACID, (1-OXOPROPOXY)-, 1-(3,3-DIMETHYLCYCLOHEXYL)-1-METH
L3
               YLETHYL ESTER"/CN
=> d 13
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
L_3
     610769-95-4 REGISTRY
RN
     Entered STN: 30 Oct 2003
ED
     Acetic acid, (1-oxopropoxy)-, 1-(3,3-dimethylcyclohexyl)-1-
CN
     methylethyl ester (9CI) (CA INDEX NAME)
FS
     3D CONCORD
MF
     C16 H28 O4
SR
     CA
                  CA, CAPLUS, CASREACT, USPATFULL
LC
     STN Files:
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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file caplus COST IN U.S. DOLLARS SINCE FILE TOTAL SESSION ENTRY FULL ESTIMATED COST 7.30 18.79 SINCE FILE TOTAL DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) ENTRY SESSION 0.00 -0.73 CA SUBSCRIBER PRICE

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=> 13

L4 1 L3

=> d l4 ti fbib abs

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN

TI Methods for the production of novel alicyclic esters having a musky smell

AN 2003:796643 CAPLUS

DN 139:307907

TI Methods for the production of novel alicyclic esters having a musky smell

IN Eh, Marcus

PA Symrise GmbH & Co. KG, Germany

SO PCT Int. Appl., 45 pp.

CODEN: PIXXD2

DT Patent

LA German

FAN. CNT 1

FAN	CNT 1																
	PATENT NO).		KIND)]	DATE		APPLICATION NO.					D?	ATE			
ΡI	WO 2003082799 WO 2003082799			A2 20031009			1	NO 2	003-1	EP329	94		20030329				
				A3													
	W: A	Æ, AG,	AL,	AM,	ΑT,	AU,	ΑZ,	BA,	BB,	BG,	BR,	BY,	ΒZ,	CA,	CH,	CN,	
	C	CO, CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	ES,	FI,	GB,	GD,	GE,	GH,	
	G	M, HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	ΚZ,	LC,	LK,	LR,	
	I	LS, LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NO,	NZ,	OM,	PH,	
	F	L, PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	ТJ,	TM,	TN,	TR,	TT,	TZ,	
	τ	JA, UG,	US,	UZ,	VC,	VN,	ΥU,	ZA,	ZM,	ZW							
		SH, GM,															
	F	(G, KZ,	MD,	RU,	TJ,	TM,	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	
	F	I, FR,	GB,	GR,	HU,	ΙE,	IT,	LU,	MC,	NL,	PT,	RO,	SE,	SI,	SK,	TR,	

BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG DE 2002-10214675 A 20020403 20031016 DE 2002-10214675 20020403 DE 10214675 A1 20030329 BR 2003004218 Α 20040727 BR 2003-4218 DE 2002-10214675 А 20020403 W 20030329 WO 2003-EP3294 20050105 20030329 EP 1492759 A2 EP 2003-732270 AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK DE 2002-10214675 A 20020403 WO 2003-EP3294 W 20030329 US 2005182273 **A1** 20050818 US 2003-510024 20030329 WO 2003-EP3294 20030329

OS CASREACT 139:307907; MARPAT 139:307907

GI

I

$$\begin{array}{c|cccc}
R^4 & R^3 & & & & & \\
R^2 & & & & & \\
R^5 & & & & \\
R^6 & & & & \\
R^7 & & & & \\
\end{array}$$
III

The invention relates to novel alicyclic esters I [R1 = Me; R2, R4 = H; R3 AB = H, Me; R5, R6 = H, Me; Y = CR7R8OC(:O)R9; R7, R8 = H, Me; R9 = HC1-5-alkyl, C2-5-alkylene; or R1, R2 = Me, Et; R3, R4 = H, Me; R5R6 = O; Y = CR7R8OC(:0)R9; or R1, R2 = Me, Et; R4, R5, R6, R7 = H, Me; Y = CR7R8OC(:0)R9], methods for their production, for their use as odorous substances for perfumed products and for odorous substance mixts. containing the inventive compds. The procedure for the preparation of I is characterized by reaction of cyclohexylalkanols II with carboxylic acids [R9CO2CR7R8CO2H, R9CO2H or XCR7R8CO2H (X = OH, halogen)] anhydrides [(R9CO2)20 or (XCR7R8CO2)20], or epoxides, III. Thus, I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2CEt] was prepared from 1-cyclohexylethanol via reaction with isobutylene oxide in cyclohexane containing BF3·OEt2, followed by reaction with (EtCO2)20 containing Et3N in the presence of catalytic DMAP. The odor of I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2CEt] was characterized (perceptible rose bloom note).

=> logoff hold SINCE FILE TOTAL COST IN U.S. DOLLARS **ENTRY** SESSION 3.10 21.89 FULL ESTIMATED COST DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL SESSION ENTRY -0.73 -1.46 CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 07:21:10 ON 29 DEC 2005